



<b>Form PTO-1449 Modified</b>		Docket No. <b>IBIS-0339</b>	Serial No. <b>09/753,869</b>
List of Patent and Publications Cited by Applicant (Use several sheets if necessary)  U.S. Department of Commerce Patent and Trademark Office		Applicant <b>Richard Griffey and Eric Swayze</b>	
		Filing Date <b>January 3, 2001</b>	Group <b>Unknown</b>
<b>OTHER DOCUMENTS (Including Author, Title, Date, Pertinent Pages, Etc.)</b>			
<i>MB</i>	<b>AA</b>	Atherton, et al., in <i>Solid Peptide Synthesis: A practical Approach</i> , IRL Press, Oxford, UK, 1989, 135	
	<b>AB</b>	Brennan, T., et al., "Two-dimensional parallel array technology as a new approach to automated combinatorial solid-phase organic synthesis," <i>Automation</i> , 1998, 33-45	
	<b>AC</b>	Chen et al., "Structure-Based Discovery of Ligands Targeted to the RNA Double Helix", <i>Biochem.</i> , 1997, Vol. 36, pp. 11402-11407	
	<b>AD</b>	Chen et al., "Spectroscopic Recognition of Guanine Dimeric Haripin Quadruplexes by a Carbocyanine Dye", <i>Proc. Natl. Acad. Sci.</i> , 1996, Vol. 93, pp. 2635-2639	
	<b>AE</b>	Corey, E.J. et al., "Computer-Assisted Analysis in Organic Synthesis," <i>Science</i> , 1985, 228, 408-418	
	<b>AF</b>	Czarnik, A.W. et al. (eds.), <i>A Practical Guide to Combinatorial Chemistry</i> , American Chemical Society, Washington, DC, 1997, Chs. 2, 13, and 14, 17-47 and 357-412	
	<b>AG</b>	Gschwend et al., "Orientational Sampling and Rigid-Body Minimization in Molecular Docking Revisited: On-the-fly Optimization and Degeneracy Removal", <i>J. Comput.-Aided Mol. Des.</i> , 1996, Vol. 10, pp. 123-132	
	<b>AH</b>	Heathcock, et al., "Total synthesis of racemic vallesamidine," <i>J. Org. Chem.</i> ®, 1990, 55(3), 798-911 (abstract only)	
	<b>AI</b>	Kuntz et al., "A Geometric Approach to Macromolecule-Ligand Interactions", <i>J. Mol. Biol.</i> , 1982, Vol. 161, pp. 269-288	
	<b>AJ</b>	Kuntz et al., "Structure-Based Molecular Design", <i>Acc. Chem. Res.</i> , 1994, Vol. 27, No. 5, pp. 117-123	
	<b>AK</b>	Levine et al., "Stalk: An Interactive System for Virtual Molecular Docking", <i>IEEE Computational Science and Engineering</i> , pp. 55-65	
	<b>AL</b>	Meng et al., "Automated Docking with Grid-Based Energy Evaluation", <i>J. Comput. Chem.</i> , 1992, Vol. 13, No. 4, pp. 505-524	
<i>MB</i>	<b>AM</b>	Rotstein, S.H. et al., "GroupBuild: A Fragment-Based Method for <i>De Novo</i> Drug Design," <i>J. Med. Chem.</i> , 1993, 36, 1700-1710	
<b>EXAMINER</b> <i>[Signature]</i>		<b>DATE CONSIDERED</b> <i>08/02</i>	





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**U. S. PATENT DOCUMENTS**

Examiner Initial		Document No.	Date	Name	Class	Subclass
<i>MM</i>	AP	5,010,175	04/23/91	Rutter, et al.	530	334
	AQ	5,186,898	02/16/93	Bridgham, et al.	422	102
	AR	5,372,672	12/13/94	Seifert, et al.	156	584
	AS	5,424,186	06/13/95	Fodor, et al.	435	006
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	AU	5,472,672	12/05/95	Brennan	422	131
	AV	5,529,756	06/25/96	Brennan	422	131
	AW	5,565,324	10/1996	Still et al.	435	6
	AX	5,573,905	11/1996	Lerner et al.	435	6
	AY	5,574,656	11/12/96	Agrafiotis et al.	364	500
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<i>MM</i>	BA	5,880,972	03/09/99	Horlbeck	364	496

**FOREIGN PATENT DOCUMENTS**

Examiner Initial		Document No.	Date	Country	Translation YES NO	

<b>EXAMINER</b> <i>MM</i>	<b>DATE CONSIDERED</b> <i>02/02</i>
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